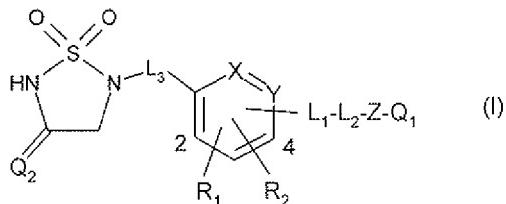


Amendment to Claims

1. (Currently Amended) A compound of the formula



wherein

R_1 is hydrogen, halogen, hydroxy, alkoxy, carboxy, cyano, nitro, trifluoromethyl, alkynyl, alkylthio, heteroaralkyl, heteroaralkoxy or heteroaryloxy provided that R_1 is located at the 2-position when L_3 is $-(CHR)_s-$ in which s is zero; or

R_1 is optionally substituted alkyl, alkenyl, optionally substituted amino, aralkyl, aralkoxy, aralkylthio, aryloxy, arylthio or cycloalkyl provided that a monocyclic aryl group which is substituted at the para position with a methylene or ethylene bridged nitrogen containing heterocycle does not constitute part of R_1 when

- (i) R_1 is located at the 2-position and L_3 is $-(CHR)_s-$ in which s is zero;
- (ii) X and Y are CH; and
- (iii) Q_2 is oxygen; or

C- R_1 may be replaced with nitrogen or N \rightarrow O; or

R_1 and R_2 combined together with the carbon atoms to which R_1 and R_2 are attached form an optionally substituted fused 5- to 6-membered aromatic or heteroaromatic ring provided that R_1 and R_2 are attached to carbon atoms adjacent to each other; or

R_2 is hydrogen, halogen, hydroxy, alkoxy, cyano, trifluoromethyl, nitro, optionally substituted amino, optionally substituted alkyl, alkylthio, aralkyl, heteroaralkyl, aralkoxy, heteroaralkoxy, aralkylthio, aryloxy, heteroaryloxy, arylthio or cycloalkyl; or

R_2 is $-C(O)R_3$ wherein

R_3 is hydroxy or optionally substituted alkoxy; or

R_3 is $-NR_4R_5$ in which R_4 and R_5 are independently hydrogen, optionally substituted alkyl, alkenyl, alkynyl, cycloalkyl, aryl, heterocyclyl, aralkyl or heteroaralkyl;

L_1 is a single bond; or

L_1 is carbon which combined together with R_2 and the carbon atoms to which L_1 and R_2 are attached form an optionally substituted fused 5- or 6-membered aromatic or heteroaromatic

ring provided that L₁ and R₂ are attached to carbon atoms adjacent to each other; or

L₁ is CH or nitrogen which taken together with R₂ and the carbon atoms to which L₁ and R₂ are attached form a fused 5- to 7-membered ring which may be interrupted with one or two heteroatoms selected from oxygen, nitrogen and sulfur provided that L₁ and R₂ are attached to carbon atoms adjacent to each other; or

L₁ is CH, oxygen, sulfur or nitrogen and L₂ is carbon which combined together with L₁, R₂ and the carbon atoms to which L₁ and R₂ are attached form an optionally substituted fused 5- or 6-membered aromatic or heteroaromatic ring provided that L₁ and R₂ are attached to carbon atoms adjacent to each other; or

L₁ is -CH₂-, oxygen, sulfur or -NR₆- and L₂ is CH which taken together with L₁, R₂ and the carbon atoms to which L₁ and R₂ are attached form a fused 5- to 7-membered ring which may be interrupted with one or two heteroatoms selected from oxygen, nitrogen and sulfur wherein

R₆ is hydrogen, optionally substituted alkyl, aralkyl, heteroaralkyl, alkoxy carbonyl, aryloxy carbonyl, carbamoyl, sulfonyl or acyl provided that L₁ and R₂ are attached to carbon atoms adjacent to each other; or

L₂ is -(CHR₇)_n- wherein

R₇ is hydrogen, hydroxy, alkoxy, carboxy, optionally substituted alkyl, cycloalkyl, aryl or heteroaryl;

n is zero or an integer from 1 to 4;

Z is -(CHR₈)_m-, -(CH₂)_mO(CHR₈)_r-, -(CH₂)_mS(CHR₈)_r- or -(CH₂)_mNR₉(CHR₈)_r- wherein

R₈ is hydrogen, optionally substituted alkyl, cycloalkyl, aryl or heterocyclyl;

R₉ is hydrogen, optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl, aralkyl, heteroaralkyl, alkoxy carbonyl, aryloxy carbonyl, heteroaryloxy carbonyl, carbamoyl, sulfonyl, acyl or acylamino;

m and r are independently zero or an integer of 1 or 2;

Q₁ is hydrogen, optionally substituted alkyl, cycloalkyl, aryl or heterocyclyl provided that

(i) Q₁ is not 2-phenyloxazol-4-yl when

R₁ and R₂ are hydrogen;

X and Y are CH;

L₁ is a single bond located at the 4-position;

L₂ is -(CHR₇)_n- wherein n is zero;

L₃ is -(CHR)_s- wherein s is zero;

Z is -(CH₂)_mO(CHR₈)_r- wherein R₈ is hydrogen, m is zero and r is 2; and

Q_2 is oxygen; or

(ii) Q_1 is not hydrogen when

R_1 and R_2 are hydrogen;

X and Y are CH ;

L_1 is a single bond;

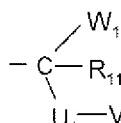
L_2 is $-(CHR_7)_n-$ wherein n is zero;

L_3 is $-(CHR)_s-$ wherein R is hydrogen and s is 1;

Z is $-(CHR_8)_m-$ wherein m is zero; and

Q_2 is oxygen; or

Q_1 is $-C(O)NR_{4a}R_{5a}$, $-C(O)R_{10}$, $-C(O)OR_{10}$ or $-S(O)_qR_{10}$ wherein R_{4a} and R_{5a} are as defined for R_4 and R_5 ; R_{10} is optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl, aralkyl or heteroaralkyl; q is an integer of 1 or 2; or



Q_1 is a radical of the formula

W_1 is aryl, heteroaryl, aralkyl or heteroaralkyl; or

W_1 is $-C(O)R_{3a}$ in which R_{3a} is hydroxy or optionally substituted alkoxy; or

R_{3a} is $-NR_{4a}R_{5a}$ in which R_{4a} and R_{5a} are as defined for R_4 and R_5 ;

R_{11} is hydrogen, alkyl or aryl;

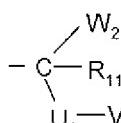
U_1 is $-C(O)-$, $-S(O)_2-$ or $-(CH_2)_r-$ in which r is as defined for Z ;

V_1 is hydroxy, alkoxy, aryl, heteroaryl, optionally substituted alkyl or cycloalkyl; or

V_1 is $-NR_{4b}R_{5b}$ in which R_{4b} and R_{5b} are as defined for R_4 and R_5 provided that

(i) L_2 is $-(CHR_7)_n-$ in which n is an integer of 1 or 2; and

(ii) Z is $-(CHR_8)_m-$ in which m is zero; or



Q_1 is a radical of the formula

W_2 is $-C(O)R_{3a}$ in which R_{3a} is hydroxy or optionally substituted alkoxy; or

R_{3a} is $-NR_{4a}R_{5a}$ in which R_{4a} and R_{5a} are as defined for R_4 and R_5 ;

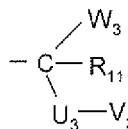
R_{11} is hydrogen, alkyl or aryl;

U_2 is $-(CH_2)_p-$ in which p is zero or 1;

V_2 is $-NR_{4b}C(O)R_{5b}$, $-NR_{4b}C(O)OR_{5b}$, $-NR_{4b}C(O)NR_{4c}R_{5b}$ or $-NR_{4b}S(O)_2R_{5b}$ in which R_{4b} and R_{4c} are as defined for R_4 , and R_{5b} has a meaning as defined for R_5 provided that

(i) L_2 is $-(\text{CH}_R)_n-$ in which n is an integer of 1 or 2; and

(ii) Z is $-(\text{CH}_R)_m-$ in which m is zero; or



Q_1 is a radical of the formula wherein

W_3 is $-\text{C}(\text{O})\text{R}_{3a}$ in which R_{3a} is hydroxy or optionally substituted alkoxy; or

R_{3a} is $-\text{NR}_{4a}\text{R}_{5a}$ in which R_{4a} and R_{5a} are as defined for R_4 and R_5 ;

R_{11} is hydrogen, alkyl or aryl;

U_3 is $-(\text{CH}_2)_p-$ in which p is zero or 1;

V_3 is $-\text{NHC}(\text{O})\text{CHR}_{4b}\text{NHC}(\text{O})\text{R}_{12}$ wherein R_{4b} is as defined for R_4 ; R_{12} is hydrogen, aryl, heterocyclyl, aralkyl, heteroaralkyl, optionally substituted alkyl, alkoxy or cycloalkyl; or

R_{12} is $-\text{NR}_{4c}\text{R}_{5b}$, in which R_{4c} and R_{5b} are as defined for R_4 and R_5 provided that

(i) L_2 is $-(\text{CH}_R)_n-$ in which n is an integer of 1 or 2; and

(ii) Z is $-(\text{CH}_R)_m-$ in which m is zero;

L_3 is $-(\text{CH}_R)_s-$ wherein

R is hydrogen, carboxy, optionally substituted alkyl, cycloalkyl, aryl or heteroaryl;

s is zero or an integer from 1 to 3;

Q_2 is oxygen, sulfur or NR_{13} wherein

R_{13} is hydrogen, hydroxy or lower alkyl;

X and Y are independently CH or nitrogen $-\text{CH}_2-$; or

~~X=Y is sulfur, oxygen or $-\text{NR}_{14}-$ wherein~~

~~R_{14} is hydrogen, optionally substituted alkyl, alkoxy carbonyl, acyl, aryloxycarbonyl, heteroaryloxycarbonyl, carbamoyl or sulfonyl;~~

or a pharmaceutically acceptable salt thereof; or a prodrug derivative thereof.

2. (Currently Amended) A The compound according to claim 1 wherein

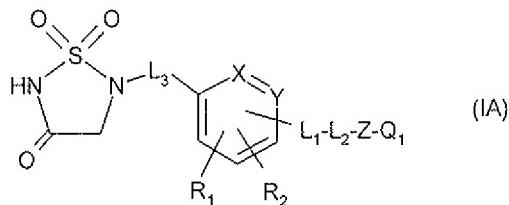
Q_2 is oxygen;

~~X and Y are CH; or~~

~~X=Y is sulfur;~~

or a pharmaceutically acceptable salt thereof; or a prodrug derivative thereof.

3. (Currently Amended) A The compound according to claim 2 of the formula



wherein

R₁ is hydrogen, halogen, hydroxy, alkoxy, trifluoromethyl, alkylthio, heteroaralkyl or heteroaralkoxy provided that R₁ is located at the 2-position when L₃ is -(CHR)_s- in which s is zero; or

R₁ is optionally substituted alkyl, aralkyl, aralkoxy or aryloxy provided that a monocyclic aryl group which is substituted at the para position with a methylene or ethylene bridged nitrogen containing heterocycle does not constitute part of R₁ when

- (i) R₁ is located at the 2-position and L₃ is -(CHR)_s- in which s is zero; and
- (ii) X and Y are CH;

R₂ is hydrogen; or

R₂ is -C(O)R₃ wherein

R₃ is hydroxy or optionally substituted alkoxy; or

R₃ is -NR₄R₅ in which R₄ and R₅ are independently hydrogen, optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl, aralkyl or heteroaralkyl;

L₁ is a single bond; or

L₁ is carbon which combined together with R₂ and the carbon atoms to which L₁ and R₂ are attached form an optionally substituted fused 5- or 6-membered aromatic or heteroaromatic ring provided that L₁ and R₂ are attached to carbon atoms adjacent to each other; or

L₁ is CH or nitrogen which taken together with R₂ and the carbon atoms to which L₁ and R₂ are attached form a fused 5- to 7-membered ring which may be interrupted with one or two heteroatoms selected from oxygen, nitrogen and sulfur provided that L₁ and R₂ are attached to carbon atoms adjacent to each other; or

L₁ is CH, oxygen, sulfur or nitrogen and L₂ is carbon which combined together with L₁, R₂ and the carbon atoms to which L₁ and R₂ are attached form an optionally substituted fused 5- or 6-membered aromatic or heteroaromatic ring provided that L₁ and R₂ are attached to carbon atoms adjacent to each other; or

L_1 is $-\text{CH}_2-$, oxygen, sulfur or $-\text{NR}_6-$ and L_2 is CH which taken together with L_1 , R_2 and the carbon atoms to which L_1 and R_2 are attached form a fused 5- to 7-membered ring which may be interrupted with one or two heteroatoms selected from oxygen, nitrogen and sulfur wherein

R_6 is hydrogen, optionally substituted alkyl, aralkyl, heteroaralkyl, alkoxy carbonyl, aryloxy carbonyl, carbamoyl, sulfonyl or acyl provided that L_1 and R_2 are attached to carbon atoms adjacent to each other; or

L_2 is $-(\text{CH}_R)_n-$ wherein

R_7 is hydrogen;

n is zero or an integer of 1 or 2;

Z is $-(\text{CHR}_8)_m-$, $-(\text{CH}_2)_m\text{O}(\text{CHR}_8)_r-$, $-(\text{CH}_2)_m\text{S}(\text{CHR}_8)_r-$ or $-(\text{CH}_2)_m\text{NR}_9(\text{CHR}_8)_r-$ wherein

R_8 is hydrogen or optionally substituted alkyl;

R_9 is hydrogen, optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl or acyl;

m and r are independently zero or an integer of 1 or 2;

Q_1 is hydrogen, optionally substituted alkyl, cycloalkyl, aryl or heterocyclyl provided that

(i) Q_1 is not 2-phenyloxazol-4-yl when

R_1 and R_2 are hydrogen;

X and Y are CH;

L_1 is a single bond located at the 4-position;

L_2 is $-(\text{CHR}_7)_n-$ wherein n is zero;

L_3 is $-(\text{CHR})_s-$ wherein s is zero; and

Z is $-(\text{CH}_2)_m\text{O}(\text{CHR}_8)_r-$ wherein R_8 is hydrogen, m is zero and r is 2; or

(ii) Q_1 is not hydrogen when

R_1 and R_2 are hydrogen;

X and Y are CH;

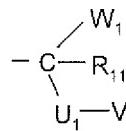
L_1 is a single bond;

L_2 is $-(\text{CHR}_7)_n-$ wherein n is zero;

L_3 is $-(\text{CHR})_s-$ wherein R is hydrogen and s is 1; and

Z is $-(\text{CHR}_8)_m-$ wherein m is zero; or

Q_1 is $-\text{C}(\text{O})\text{NR}_{4a}\text{R}_{5a}$, $-\text{C}(\text{O})\text{R}_{10}$, $-\text{C}(\text{O})\text{OR}_{10}$ or $-\text{S}(\text{O})_q\text{R}_{10}$ wherein R_{4a} and R_{5a} are as defined for R_4 and R_5 ; R_{10} is optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl, aralkyl or heteroaralkyl; q is an integer of 1 or 2; or



Q_1 is a radical of the formula $-\begin{array}{c} W_1 \\ \diagdown \\ C - R_{11} \\ \diagup \\ U_1 - V_1 \end{array}$ wherein

W_1 is aryl, heteroaryl, aralkyl or heteroaralkyl; or

W_1 is $-C(O)R_{3a}$ in which R_{3a} is hydroxy or optionally substituted alkoxy; or

R_{3a} is $-NR_{4a}R_{5a}$ in which R_{4a} and R_{5a} are as defined for R_4 and R_5 ;

R_{11} is hydrogen, alkyl or aryl;

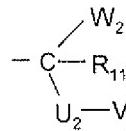
U_1 is $-C(O)-$ or $-(CH_2)_r-$ in which r is as defined for Z ;

V_1 is hydroxy, alkoxy, aryl, heteroaryl, optionally substituted alkyl or cycloalkyl; or

V_1 is $-NR_{4b}R_{5b}$ in which R_{4b} and R_{5b} are as defined for R_4 and R_5 provided that

(i) L_2 is $-(CHR_7)_n-$ in which n is an integer of 1 or 2; and

(ii) Z is $-(CHR_8)_m-$ in which m is zero; or



Q_1 is a radical of the formula $-\begin{array}{c} W_2 \\ \diagdown \\ C - R_{11} \\ \diagup \\ U_2 - V_2 \end{array}$ wherein

W_2 is $-C(O)R_{3a}$ in which R_{3a} is hydroxy or optionally substituted alkoxy; or

R_{3a} is $-NR_{4a}R_{5a}$ in which R_{4a} and R_{5a} are as defined for R_4 and R_5 ;

R_{11} is hydrogen, alkyl or aryl;

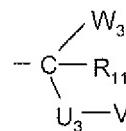
U_2 is $-(CH_2)_p-$ in which p is zero or 1;

V_2 is $-NR_{4b}C(O)R_{5b}$, $-NR_{4b}C(O)OR_{5b}$, $-NR_{4b}C(O)NR_{4c}R_{5b}$ or $-NR_{4b}S(O)_2R_{5b}$ in which

R_{4b} and R_{4c} are as defined for R_4 , and R_{5b} has a meaning as defined for R_5 provided that

(i) L_2 is $-(CHR_7)_n-$ in which n is an integer of 1 or 2; and

(ii) Z is $-(CHR_8)_m-$ in which m is zero; or



Q_1 is a radical of the formula $-\begin{array}{c} W_3 \\ \diagdown \\ C - R_{11} \\ \diagup \\ U_3 - V_3 \end{array}$ wherein

W_3 is $-C(O)R_{3a}$ in which R_{3a} is hydroxy or optionally substituted alkoxy; or

R_{3a} is $-NR_{4a}R_{5a}$ in which R_{4a} and R_{5a} are as defined for R_4 and R_5 ;

R_{11} is hydrogen, alkyl or aryl;

U_3 is $-(CH_2)_p-$ in which p is zero or 1;

V_3 is $-NHC(O)CHR_{4b}NHC(O)R_{12}$ wherein R_{4b} is as defined for R_4 ; R_{12} is hydrogen, aryl, heterocyclyl, aralkyl, heteroaralkyl, optionally substituted alkyl, alkoxy or cycloalkyl; or

R_{12} is $-NR_{4c}R_{5b}$, in which R_{4c} and R_{5b} are as defined for R_4 and R_5 provided that

- (i) L_2 is $-(CH_7)_n-$ in which n is an integer of 1 or 2; and
- (ii) Z is $-(CHR_8)_m-$ in which m is zero;

L_3 is $-(CHR)_s-$ wherein

R is hydrogen;

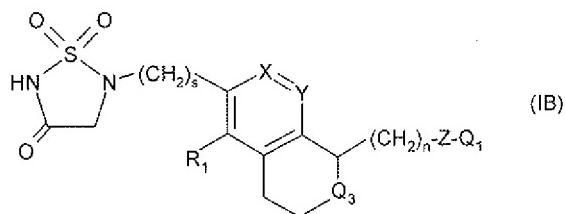
s is zero or an integer from 1 to 3;

X and Y are CH; or

~~X=Y is sulfur;~~

or a pharmaceutically acceptable salt thereof; or a prodrug derivative thereof.

4. (Currently Amended) A The compound according to claim 3 of the formula



wherein

R_1 is hydrogen, halogen, hydroxy, alkoxy, trifluoromethyl, optionally substituted alkyl, alkylthio, aralkyl, aralkoxy, aryloxy, heteroaralkyl or heteroaralkoxy;

n is zero or an integer of 1 or 2;

Z is $-(CHR_8)_m-$, $-(CH_2)_mO(CHR_8)_r-$, $-(CH_2)_mS(CHR_8)_r-$ or $-(CH_2)_mNR_9(CHR_8)_r-$ wherein

R_8 is hydrogen;

R_9 is hydrogen, optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl or acyl;

m and r are independently zero or an integer of 1 or 2;

Q_1 is hydrogen, optionally substituted alkyl, cycloalkyl, aryl or heterocyclyl; or

Q_1 is $-C(O)NR_{4a}R_{5a}$, $-C(O)R_{10}$, $-C(O)OR_{10}$ or $-S(O)_qR_{10}$ wherein

R_{4a} and R_{5b} are independently hydrogen, optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl, aralkyl or heteroaralkyl;

R_{10} is optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl, aralkyl or heteroaralkyl;

q is an integer of 1 or 2;

s is zero or an integer of 1 or 2;

Q₃ is O, S or -NR_{6a-} wherein

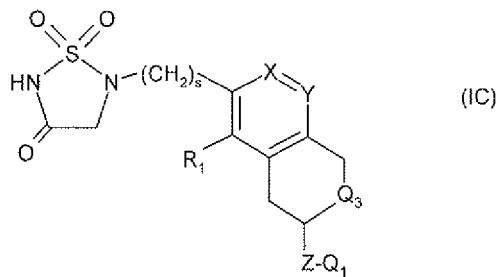
R_{6a} is hydrogen, optionally substituted alkyl, aralkyl, heteroaralkyl, alkoxy carbonyl, aryloxy carbonyl, carbamoyl, sulfonyl or acyl;

X and Y are CH; or

~~X=Y is sulfur;~~

or a pharmaceutically acceptable salt thereof; or a prodrug derivative thereof.

5. (Currently Amended) A The compound according to claim 3 of the formula



wherein

R₁ is hydrogen, halogen, hydroxy, alkoxy, trifluoromethyl, optionally substituted alkyl, alkylthio, aralkyl, aralkoxy, aryloxy, heteroaralkyl or heteroaralkoxy;

Z is -(CHR₈)_m- , -(CH₂)_mO(CHR₈)_r- , -(CH₂)_mS(CHR₈)_r- or -(CH₂)_mNR₉(CHR₈)_r- wherein

R₈ is hydrogen;

R₉ is hydrogen, optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl or acyl;

m and r are independently zero or an integer of 1 or 2;

Q₁ is hydrogen, optionally substituted alkyl, cycloalkyl, aryl or heterocyclyl; or

Q₁ is -C(O)NR_{4a}R_{5a}, -C(O)R₁₀, -C(O)OR₁₀ or -S(O)_qR₁₀ wherein

R_{4a} and R_{5a} are independently hydrogen, optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl, aralkyl or heteroaralkyl;

R₁₀ is optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl, aralkyl or heteroaralkyl;

q is an integer of 1 or 2;

s is zero or an integer of 1 or 2;

Q₃ is O, S or -NR_{6a-} wherein

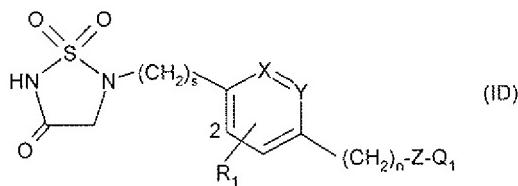
R_{6a} is hydrogen, optionally substituted alkyl, aralkyl, heteroaralkyl, alkoxy carbonyl,

aryloxycarbonyl, carbamoyl, sulfonyl or acyl;
X and Y are CH; or
~~-X=Y-~~ is sulfur;
or a pharmaceutically acceptable salt thereof; or a prodrug derivative thereof.

6. (Currently Amended) A The compound according to claim 3 wherein

R₂ is hydrogen;
L₁ is a single bond;
L₂ is -(CH₂)_n- in which n is zero or an integer of 1 or 2;
or a pharmaceutically acceptable salt thereof; or a prodrug derivative thereof.

7. (Currently Amended) A The compound according to claim 6 of the formula



wherein

R₁ is hydrogen, halogen, hydroxy, alkoxy, trifluoromethyl or alkylthio provided that R₁ is located at the 2-position when s is zero; or

R₁ is optionally substituted alkyl, aralkyl, aralkoxy or aryloxy provided that a monocyclic aryl group which is substituted at the para position with a methylene or ethylene bridged nitrogen containing heterocycle does not constitute part of R₁ when

- (i) R₁ is located at the 2-position and s is zero; and
- (ii) X and Y are CH;

n is zero or an integer of 1 or 2;

s is zero or 1;

Z is -(CHR₈)_m-, -(CH₂)_mO(CHR₈)_r-, -(CH₂)_mS(CHR₈)_r- or -(CH₂)_mNR₉(CHR₈)_r- wherein

R₈ is hydrogen;

R₉ is hydrogen, optionally substituted alkyl, cycloalkyl, aryl, heteroaryl or acyl;

m and r are independently zero or an integer of 1 or 2;

Q_1 is hydrogen, optionally substituted alkyl, cycloalkyl, aryl or heterocyclyl provided that

(i) Q_1 is not 2-phenyloxazol-4-yl when

R_1 is hydrogen;

X and Y are CH;

n is zero;

s is zero; and

Z is $-(CH_2)_mO(CHR_8)_r-$ wherein R_8 is hydrogen, m is zero and r is 2; or

(ii) Q_1 is not hydrogen when

R_1 is hydrogen;

X and Y are CH;

n is zero;

s is 1;

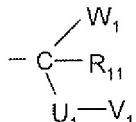
Z is $-(CHR_8)_m-$ wherein m is zero; or

Q_1 is $-C(O)NR_{4a}R_{5a}$, $-C(O)R_{10}$, $-C(O)OR_{10}$ or $-S(O)_qR_{10}$ wherein

R_{4a} and R_{5a} are independently hydrogen, optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl, aralkyl or heteroaralkyl;

R_{10} is optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl, aralkyl or heteroaralkyl;

q is an integer of 1 or 2; or



Q_1 is a radical of the formula

W_1 is aryl, heteroaryl, aralkyl or heteroaralkyl; or

W_1 is $-C(O)R_{3a}$ in which R_{3a} is hydroxy or optionally substituted alkoxy; or

R_{3a} is $-NR_{4a}R_{5a}$ in which R_{4a} and R_{5a} are independently hydrogen, optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl, aralkyl or heteroaralkyl;

R_{11} is hydrogen, alkyl or aryl;

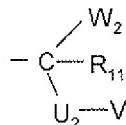
U_1 is $-C(O)-$ or $-(CH_2)_r-$ in which r is as defined for Z;

V_1 is hydroxy, alkoxy, aryl, heteroaryl, optionally substituted alkyl or cycloalkyl; or

V_1 is $-NR_{4b}R_{5b}$ in which R_{4b} and R_{5b} are as defined for R_{4a} and R_{5a} provided that

(i) n is an integer of 1 or 2; and

(ii) Z is $-(CHR_8)_m-$ in which m is zero; or



Q_1 is a radical of the formula $-\begin{array}{c} \diagup \\ W_2 \\ \diagdown \\ C \\ \diagup \\ R_{11} \\ \diagdown \\ U_2 \end{array} - V_2$ wherein

W_2 is $-C(O)R_{3a}$ in which R_{3a} is hydroxy or optionally substituted alkoxy; or

R_{3a} is $-NR_{4a}R_{5a}$ in which R_{4a} and R_{5a} are independently hydrogen, optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl, aralkyl or heteroaralkyl;

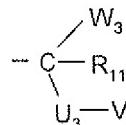
R_{11} is hydrogen, alkyl or aryl;

U_2 is $-(CH_2)_p-$ in which p is zero or 1;

V_2 is $-NR_{4b}C(O)R_{5b}$, $-NR_{4b}C(O)OR_{5b}$, $-NR_{4b}C(O)NR_{4c}R_{5b}$ or $-NR_{4b}S(O)_2R_{5b}$ in which R_{4b} and R_{4c} are as defined for R_{4a} , and R_{5b} has a meaning as defined for R_{5a} provided that

(i) n is an integer of 1 or 2; and

(ii) Z is $-(CHR_8)_m-$ in which m is zero; or



Q_1 is a radical of the formula $-\begin{array}{c} \diagup \\ W_3 \\ \diagdown \\ C \\ \diagup \\ R_{11} \\ \diagdown \\ U_3 \end{array} - V_3$ wherein

W_3 is $-C(O)R_{3a}$ in which R_{3a} is hydroxy or optionally substituted alkoxy; or

R_{3a} is $-NR_{4a}R_{5a}$ in which R_{4a} and R_{5a} are independently hydrogen, optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl, aralkyl or heteroaralkyl;

R_{11} is hydrogen, alkyl or aryl;

U_3 is $-(CH_2)_r-$ in which r is zero or 1;

V_3 is $-NHC(O)CHR_{4b}NHC(O)R_{12}$ wherein R_{4b} is as defined for R_{4a} ; R_{12} is hydrogen, aryl, heterocyclyl, aralkyl, heteroaralkyl, optionally substituted alkyl, alkoxy or cycloalkyl; or

R_{12} is $-NR_{4c}R_{5b}$ in which R_{4c} is as defined for R_{4a} , and R_{5b} has a meaning as defined for R_{5a} provided that

(i) n is an integer of 1 or 2; and

(ii) Z is $-(CHR_8)_m-$ in which m is zero;

X and Y are CH; or

~~X=Y is sulfur;~~

or a pharmaceutically acceptable salt thereof; or a prodrug derivative thereof.

8. (Canceled).

9. (Currently Amended) A The compound according to claim 7 wherein

R₁ is bromide;

X and Y are CH;

or a pharmaceutically acceptable salt thereof; or a prodrug derivative thereof.

10. (Currently Amended) A The compound according to claim 7 wherein

n is zero;

s is 1;

Z is -(CH₂)_m- in which m is zero;

Q₁ is -C(O)NR_{4a}R_{5a}, -C(O)R₁₀, -C(O)OR₁₀ or -S(O)_qR₁₀ wherein

R_{4a} and R_{5a} are independently hydrogen, optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl, aralkyl or heteroaralkyl;

R₁₀ is optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl, aralkyl or heteroaralkyl;

q is an integer of 1 or 2;

or a pharmaceutically acceptable salt thereof; or a prodrug derivative thereof.

11. (Currently Amended) A The compound according to claim 7 wherein

n is an integer of 1 or 2;

Z is -(CH₂)_mO(CH₂)_r- or -(CH₂)_mS(CH₂)_r- wherein

m is zero;

r is zero or 1;

Q₁ is optionally substituted alkyl, cycloalkyl, aryl or heterocyclyl;

or a pharmaceutically acceptable salt thereof; or a prodrug derivative thereof.

12. (Currently Amended) A The compound according to claim 7 wherein

n is an integer of 1 or 2;

Z is -(CH₂)_mNR₉(CH₂)_r- wherein

R₉ is hydrogen, optionally substituted alkyl, cycloalkyl, aryl, heteroaryl or acyl;

m is zero;

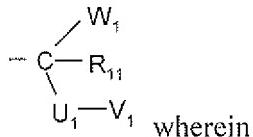
r is zero or 1;

Q_1 is optionally substituted alkyl, cycloalkyl, aryl or heterocyclyl; or
 Q_1 is $-C(O)NR_{4a}R_{5a}$, $-C(O)R_{10}$, $-C(O)OR_{10}$ or $-S(O)_qR_{10}$ wherein
 R_{4a} and R_{5a} are independently hydrogen, optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl, aralkyl or heteroaralkyl;
 R_{10} is optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl, aralkyl or heteroaralkyl;
 q is an integer of 1 or 2;
or a pharmaceutically acceptable salt thereof; or a prodrug derivative thereof.

13. (Currently Amended) A The compound according to claim 7 wherein

n is an integer of 1 or 2;

Z is $-(CH_2)_m-$ wherein m is zero;



Q_1 is a radical of the formula

W_1 is aryl, heteroaryl, aralkyl or heteroaralkyl;

R_{11} is hydrogen, alkyl or aryl;

U_1 is $-C(O)-$ or $-(CH_2)_r-$ in which r is zero;

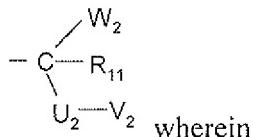
V_1 is aryl, heteroaryl, optionally substituted alkyl or cycloalkyl;

or a pharmaceutically acceptable salt thereof; or a prodrug derivative thereof.

14. (Currently Amended) A The compound according to claim 7 wherein

n is 1;

Z is $-(CH_2)_m-$ wherein m is zero;



Q_1 is a radical of the formula

W_2 is $-C(O)R_{3a}$ in which R_{3a} is $-NR_{4a}R_{5a}$, and R_{4a} and R_{5a} are independently hydrogen, optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl, aralkyl or heteroaralkyl;

R_{11} is hydrogen;

U_2 is $-(CH_2)_p-$ in which p is zero;

V_2 is $-NR_{4b}C(O)R_{5b}$, $-NR_{4b}C(O)OR_{5b}$, $-NR_{4b}C(O)NR_{4c}R_{5b}$ or $-NR_{4b}S(O)_2R_{5b}$ in which

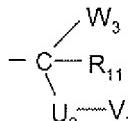
R_{4b} and R_{4c} are as defined for R_{4a} , and R_{5b} has a meaning as defined for R_{5a} ;

or a pharmaceutically acceptable salt thereof; or a prodrug derivative thereof.

15. (Currently Amended) A The compound according to claim 7 wherein

n is 1;

Z is -(CH₂)_m- wherein m is zero;



Q₁ is a radical of the formula

W₃ is -C(O)R_{3a} in which R_{3a} is -NR_{4a}R_{5a}, and R_{4a} and R_{5a} are independently hydrogen, optionally substituted alkyl, cycloalkyl, aryl, heterocyclyl, aralkyl or heteroaralkyl;

R₁₁ is hydrogen;

U₃ is -(CH₂)_p- in which p is zero;

V₃ is -NHC(O)CHR_{4b}NHC(O)R₁₂ wherein R_{4b} is as defined for R_{4a}; R₁₂ is hydrogen, aryl, heterocyclyl, aralkyl, heteroaralkyl, optionally substituted alkyl or alkoxy; or

R₁₂ is -NR_{4c}R_{5b} in which R_{4c} and R_{5b} are as defined for R_{4a} and R_{5a};

or a pharmaceutically acceptable salt thereof; or a prodrug derivative thereof.

16. (Currently Amended) A The compound according to claim 1 which is selected from:

5-Naphthalen-1-ylmethyl-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

N-[3-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzyl]-acetamide;

[3-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzyl]-carbamic acid *t*-butyl ester;

5-(4-Aminomethyl-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

N-[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzyl]-acetamide;

[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzyl]-carbamic acid *t*-butyl ester;

3-Phenyl-N-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzyl]-propionamide;

5-(3-Iodo-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

5-(3-Nitro-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

5-(3-Amino-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

N-[3-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenyl]-acetamide;

4,1-Dioxo-5-pyridin-4-ylmethyl-1,2,5-thiadiazolidin-3-one;

5-(4-Amino-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

N-[3-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenyl]-butyramide;

1-Propyl-3-[3-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenyl]-urea;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid methyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid;

2-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid;

5-(2-Methyl-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

~~4,1-Dioxo-5-pyridin-3-ylmethyl-1,2,5-thiadiazolidin-3-one;~~
~~4,1-Dioxo-5-pyridin-2-ylmethyl-1,2,5-thiadiazolidin-3-one;~~
~~5-(6-Amino-pyridin-3-ylmethyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;~~
~~4,1-Dioxo-5-thiophen-2-ylmethyl-1,2,5-thiadiazolidin-3-one;~~
~~5-(4-Methoxy-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;~~
~~5-(4-Amino-2-bromo-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;~~
~~N-[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenyl]-acetamide;~~
~~N-[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenyl]-methanesulfonamide;~~
~~N-[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzyl]-methanesulfonamide;~~
~~5-(4-Methyl-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;~~
~~Amino-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenyl]-acetic acid;~~
~~2-Amino-N-propyl-2-[2-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenyl]-acetamide;~~
~~2-Amino-N-propyl-2-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenyl]-acetamide;~~
2,2,2-Trifluoro-N-{propylcarbamoyl-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenyl]-methyl}-acetamide;
2-Methanesulfonylamino-N-propyl-2-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenyl]-acetamide;
2-Acetylamino-N-propyl-3-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenyl]-propionamide;
2-Acetylamino-2-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzyl]-malonic acid diethyl ester;
2-Amino-N-propyl-3-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenyl]-propionamide;
2-Acetylamino-3-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenyl]-propionic acid ethyl ester;
Phenyl-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-acetic acid;
1,1-Dioxo-5-phenethyl-1,2,5-thiadiazolidin-3-one;
5-[2-(4-Methyl-thiazol-5-yl)-ethyl]-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-[2-(3,4-Dimethoxy-phenyl)-ethyl]-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-[2-(2-Chloro-phenyl)-ethyl]-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-[2-(4-Amino-phenyl)-ethyl]-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
2,2,2-Trifluoro-N-{4-[2-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-ethyl]-phenyl}-acetamide;

N-{4-[2-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-yl)-ethyl]-phenyl}-butyramide;
4,1-Dioxo-5-(2-pyridin-3-yl-ethyl)-1,2,5-thiadiazolidin-3-one;
4,1-Dioxo-5-(2-pyridin-4-yl-ethyl)-1,2,5-thiadiazolidin-3-one;
3-Phenyl-2-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-propionic acid;
5-[2-(3-Amino-phenyl)-ethyl]-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-(4-Aminomethyl-naphthalen-1-ylmethyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-(1-Ethyl-2-methyl-1H-benzimidazol-5-ylmethyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-[2-Methyl-1-(3-methyl-butyl)-1H-benzimidazol-5-ylmethyl]-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-(4-Methoxy-quinolin-7-ylmethyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-(4-Isobutoxy-quinolin-7-ylmethyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
{(1-Butylcarbamoyl-3-phenyl-propyl)-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoyl]-amino}-acetic acid;
{[Butylcarbamoyl-(4-ethyl-phenyl)-methyl]-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoyl]-amino}-acetic acid;
{[Butylcarbamoyl-(3-phenoxy-phenyl)-methyl]-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoyl]-amino}-acetic acid;
{[Butylcarbamoyl-(4-methoxy-phenyl)-methyl]-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoyl]-amino}-acetic acid;
{[(2-Bromo-phenyl)-butylcarbamoyl-methyl]-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoyl]-amino}-acetic acid;
(Butylcarbamoyl-naphthalen-2-yl-methyl)-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoyl]-amino}-acetic acid;
{[Butylcarbamoyl-(4-chloro-phenyl)-methyl]-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoyl]-amino}-acetic acid;
{[(3-Benzylxy-phenyl)-butylcarbamoyl-methyl]-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoyl]-amino}-acetic acid;
{((E)-1-Butylcarbamoyl-3-phenyl-allyl)-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoyl]-amino}-acetic acid;
N-(1-Butylcarbamoyl-3-phenyl-propyl)-N-(4-(1,1,4-trioxo-1,2,5-thiazodiazolidin-2-ylmethyl)-benzoyl)-amino-acetic acid;
4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 4-methanesulfonyl-benzyl ester;
4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 3-chloro-benzyl ester;
4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 4-butyl-benzyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 4-hydroxymethyl-benzyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 2-phenethyl-benzyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid biphenyl-2-ylmethyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 4-difluoromethoxy-benzyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 5-(carboxy-difluoro-methyl)-thiophen-2-ylmethyl ester;

[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenylmethanesulfonyl]-acetic acid ethyl ester;

[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzylsulfanyl]-acetic acid ethyl ester;

5-[4-(3-Methyl-butylsulfanyl-methyl)-benzyl]-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 2-ethyl-butyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid cyclobutylmethyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid cyclopentylmethyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 2-methyl-pentyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 2,4,4-trimethyl-pentyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid cyclohexylmethyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 1,2-dimethyl-propyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid cyclopentyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 2-methyl-butyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 2-methylsulfanyl-ethyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 2-carboxymethylsulfanyl-ethyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 5-nitro-furan-2-ylmethyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid pyridin-2-ylmethyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 3-hydroxymethyl-benzyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 3-methanesulfonyl-benzyl ester;

(4-{4-[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoylamino]-butyl}-phenyl)-acetic acid;

(4-{3-[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoylamino]-propyl}-phenyl)-acetic acid;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 5-dimethylaminomethyl-furan-2-ylmethyl ester;

(S)-2-Acetylamino-N-[(S)-1-pentylcarbamoyl-2-[3-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenyl]-ethyl]-3-phenyl-propionamide;

5-(1H-Indol-5-ylmethyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

1,1-Dioxo-5-(3,4,5-trimethoxy-benzyl)-1,2,5-thiadiazolidin-3-one;

5-[4-(4-Benzyl-piperazin-1-ylmethyl)-benzyl]-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenyl]-acetic acid;

5-(4-Benzoyl-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

5-Naphthalen-2-ylmethyl-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

5-[4-(4-Methyl-pentanoyl)-benzyl]-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

5-[3-(2-Fluoro-phenoxy)-benzyl]-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

3-{2-[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenyl]-ethoxy}-benzoic acid;

1-(3-Methyl-butyl)-6-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-1H-quinolin-2-one;

5-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-thiophene-2-carboxylic acid methyl phenethyl amide;

5-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-thiophene-2-carboxylic acid (2-thiophen-2-yl-ethyl) amide;

5-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-thiophene-2-carboxylic acid phenethyl amide;

[4-(2-{[5-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-thiophene-2-carbonyl] amino}-ethyl)-phenyl]-acetic acid;

5-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-thiophene-2-carboxylic acid 4-carboxy-benzyl ester;

5-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-thiophene-2-carboxylic acid isobutyl ester;

5-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-thiophene-2-carboxylic acid isobutyl amide;

2-Amino-N-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzyl]-acetamide;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 4-carboxy-benzyl ester;

1,1-Dioxo-5-(3-phenoxy-benzyl)-1,2,5-thiadiazolidin-3-one;

3-Nitro-4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid;

5-(4-Hydroxymethyl-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

2-Amino-4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid methyl ester;
5-(4-Hydroxy-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-Nitro-2-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid;
5-Amino-2-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid;
5-(4-Chloro-3-methoxy-5-nitro-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-(2-Nitro-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-(3-Methyl-2-nitro-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-(3-Methyl-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
1,1-Dioxo-5-(3-phenyl-propyl)-1,2,5-thiadiazolidin-3-one;
5-(4-Butoxy-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
1,1-Dioxo-5-(2-trifluoromethyl-benzyl)-1,2,5-thiadiazolidin-3-one;
3-Amino-4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid;
4-[5-Amino-2-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenyl]-butyric acid;
5-(2-Methyl-3-nitro-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-(4-Methyl-3-nitro-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-(5-Methyl-2-nitro-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-(2-Amino-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
2-[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzyl]-isoindole-1,3-dione;
2-[3-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzyl]-isoindole-1,3-dione;
5,5'-[1,4-Phenylenebis(methylene)bis[1,2,5-thiadiazolidine-3-one], 1,1-dioxide;
N-[2-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenyl]-oxalamic acid;
5-(3-Hydroxy-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
2-Amino-4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid;
5-[5-(4-Nitro-phenyl)-furan-2-ylmethyl]-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-(4-Fluoro-2-trifluoromethyl-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-(3-Hydroxymethyl-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-(3-Amino-5-hydroxymethyl-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-(3-Amino-4-methyl-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-(2-Amino-3-methyl-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-(3-Amino-2-methyl-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-(2-Amino-5-methyl-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
2,2,2-Trifluoro-N-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzyl]-acetamide;
4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-pyridine-2-carbonitrile;
4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-pyridine-2-carboxylic acid ethyl ester;
5-(3,4-Dimethoxy-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

5-(3-Amino-5-hydroxy-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-(3,5-Dimethyl-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
(S)-3-Phenyl-2-[3-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzylamino]-propionic acid ethyl ester;
(S)-3-Phenyl-2-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzylamino]-propionic acid ethyl ester;
2-Amino-5-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid methyl ester;
2-Acetylamino-4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid methyl ester;
5-(2-Benzyl-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5-(2,4-Bis-trifluoromethyl-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
1,1-Dioxo-5-(2,4,6-trifluoro-benzyl)-1,2,5-thiadiazolidin-3-one;
5-(2-Bromo-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
5,5'-[[1,1'-biphenyl]-2,2'-diyl]bis(methylene)bis[1,2,5-thiadiazolidine-3-one], 1,1-dioxide;
5-(4-Ethylaminomethyl-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
2-Acetylamino-4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid;
2-Amino-4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid ethyl ester;
1,1-Dioxo-5-[4-(phenethylamino-methyl)-benzyl]-1,2,5-thiadiazolidin-3-one;
5-(4-Diethylaminomethyl-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
2-Amino-4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid benzyl ester;
N-Benzyl-4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzamide;
5-(5-Dimethylaminomethyl-furan-2-ylmethyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;
N-[2-(3-Trifluoromethyl-phenyl)-ethyl]-4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzamide;
N-(3-Methyl-butyl)-4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzamide;
(S)-3-Phenyl-2-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-propionic acid;
(R)-3-Phenyl-2-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-propionic acid;
4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid benzyl ester;
[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenoxy]-acetic acid;
4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid isobutyl ester;
2-Amino-4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid isobutyl ester;
[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenoxy]-acetic acid methyl ester;
4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 4-carboxymethoxy-benzyl ester;
5-(5-Aminomethyl-thiophen-2-ylmethyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

4-{2-[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzylamino]-ethyl}-benzoic acid;

[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenoxy]-acetic acid isobutyl ester;

[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-phenoxy]-acetic acid benzyl ester;

N-Isobutyl-4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzamide;

5-(5-Diethylaminomethyl-thiophen-2-ylmethyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

4-(2-{{[5-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-thiophen-2-ylmethyl]-amino}-ethyl}-benzoic acid;

3-Nitro-4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid methyl ester;

3-Nitro-4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid ethyl ester;

3-Nitro-4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid isobutyl ester;

5-(4-Ethoxy-benzyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

1,1-Dioxo-5-(3-trifluoromethyl-benzyl)-1,2,5-thiadiazolidin-3-one;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 4-carboxymethyl-benzyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid phenethyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 2-phenylamino-ethyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 2-(3-methoxy-phenyl)-ethyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 2,2-dimethyl-propyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 2-methoxycarbonyl-2-methyl-propyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 2,2,4-trimethyl-pentyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 3-dimethylamino-2,2-dimethyl-propyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid (3aR,4S,5R,6aS)-5-benzoyloxy-2-oxo-hexahydro-cyclopenta[b]furan-4-ylmethyl ester;

6-{{[5-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-thiophen-2-ylmethyl]-amino}-hexanoic acid;

5-[(3-Methyl-butylamino)-methyl]-thiophen-2-ylmethyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 3-methyl-4-nitro-benzyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 3-chloro-4-methyl-benzyl ester;

5-[5-(Isobutylamino-methyl)-thiophen-2-ylmethyl]-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 5-ethoxycarbonyl-pentyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 2-(3-chloro-phenyl)-ethyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 2-m-tolyl-ethyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 2-(3-trifluoromethyl-phenyl)-ethyl ester;

(R)-3-Phenyl-2-[4-(1,1,4-triexo-1,2,5-thiadiazolidin-2-ylmethyl)-benzylamino]-propionic acid ethyl ester;

5-[4-(Benzylamino-methyl)-benzyl]-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 4-methyl-benzyl ester;

4-Methyl-6-{{[5-(1,1,4-triexo-1,2,5-thiadiazolidin-2-ylmethyl)-thiophen-2-ylmethyl]}-amino}-hexanoic acid;

4-[(1,1,4-tribedo-1,2,5-thiadiazolidin-2-yl)methyl]-benzoic acid [4-(methoxycarbonyl)-phenyl]methyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 2-cyclohexyl-2-methyl-propyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 2-phenoxy-propyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 4-trifluoromethyl-benzyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 3-trifluoromethyl-benzyl ester;

4-[(1,1,4-tribedo-1,2,5-thiadiazolidin-2-yl)methyl]-benzoic acid 2-(4-carboxyphenyl)ethyl ester;

5-[5-(3-Methylbutyryl)-thiophen-2-ylmethyl]-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

3-[[4-[(1,1,4-Tribedo-1,2,5-thiadiazolidin-2-yl)methyl]benzoyl]-oxy]methyl]benzoic acid;

5-[4-(Isobutylamino-methyl)-benzyl]-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

5-{4-[(2,2-Dimethyl-propylamino)-methyl]-benzyl}-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid naphthalen-1-ylmethyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 4-nitro-benzyl ester;

(4-{2-[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoylamino]-ethyl}-phenyl)-acetic acid;

~~5-[5-(4-Methyl-pentanoyl)-thiophen-2-ylmethyl]-1,1-dioxo-1,2,5-thiadiazolidin-3-one;~~

~~5-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-thiophene-2-carboxylic acid;~~

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 3-nitro-benzyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 3-(carboxymethyl-amino)-2,2-dimethyl-propyl ester;

5-[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoyloxymethyl]-thiophene-2-carboxylic acid;

5-[4-(4-Benzyl-piperazin-1-ylmethyl)-benzyl]-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid biphenyl-4-ylmethyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 4-acetylamino-benzyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 2-benzyl-benzyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 2-methyl-benzyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 2-methyl-3-nitro-benzyl ester;

Glycine, N-(aminosulfonyl)-N-[[4-[[[(2-phenylethyl)thio]methyl]phenyl]methyl]-, methyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 3-carboxymethyl-benzyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 4-methyl-3-nitro-benzyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 4-fluoro-2-trifluoromethyl-benzyl ester;

4-[5-(2,4-Dimethoxy-benzyl)-1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl]-benzoic acid 4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-benzyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-benzyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 5-methyl-2-nitro-benzyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid o-tolyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 3-(carboxymethyl-methyl-amino)-2,2-dimethyl-propyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid phenyl ester

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 5-isobutylcarbamoyl-thiophen-2-ylmethyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid naphthalen-2-ylmethyl ester;

N,N-Diisobutyl-4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzamide;

{4-[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoyl]-piperazin-1-yl}-acetic acid;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid naphthalen-2-yl ester;

5-[4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoyloxymethyl]-thiophene-2-carboxylic acid isobutyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 5-carbamoyl-thiophen-2-ylmethyl ester;

5-[4-(4-Benzyl-piperazine-1-carbonyl)-benzyl]-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 5-(3-phenyl-propionyl)-thiophen-2-ylmethyl ester;

4-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-ylmethyl)-benzoic acid 5-benzylcarbamoyl-thiophen-2-ylmethyl ester;

1,1-Dioxo-5-phenyl-1,2,5-thiadiazolidin-3-one;

5-(2,4-Diamino-phenyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

3-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-yl)-benzoic acid methyl ester;

3-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-yl)-benzoic acid;

5-(4-Aminomethyl-phenyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

[2-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-yl)-phenyl]-acetic acid methyl ester;

[2-(1,1,4-Trioxo-1,2,5-thiadiazolidin-2-yl)-phenyl]-acetic acid;

5-(2,4-Dimethoxyphenyl)-1,1-dioxo-[1,2,5]thiadiazolidin-3-one potassium salt;

N-Benzyl-2-[3-methyl-4-(1,1,4-trioxo-[1,2,5]thiadiazolidin-2-yl)-phenoxy]-acetamide;

3-[3-Hydroxy-4-(1,1,4-trioxo-[1,2,5]thiadiazolidin-2-yl)-benzyl]-3,4-dihydro-1H-benzo[1,4]diazepine-2,5-dione;

5-(4-Iodo-phenyl)-1,1-dioxo-1,2,5-thiadiazolidin-3-one;

(S)-2-Amino-3-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-phenyl]-propionic acid benzyl ester;

(S)-2-Amino-3-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-phenyl]-propionic acid;

(S)-2-Acetylamino-N-[(S)-1-pentylcarbamoyl-2-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-phenyl]-ethyl]-3-phenyl-propionamide;

(S)-2-Acetylamino-3-phenyl-N-[(S)-1-(4-phenyl-butylcarbamoyl)-2-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-phenyl]-ethyl]-propionamide;

[4-(2-[(S)-2-((S)-2-Acetylamino-3-phenyl-propionylamino)-3-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-phenyl]-propionylamino}-ethyl]-phenyl]-acetic acid;

2-[4-(2-Benzoylamino-2-{1-carbamoyl-2-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-phenyl]-ethylcarbamoyl}-ethyl)-phenoxy]-malonic acid;

(S)-2-(Biphenyl-4-sulfonylamino)-N-pentyl-3-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-phenyl]-propionamide;

(S)-2-(Biphenyl-4-sulfonylamino)-N-(4-phenyl-butyl)-3-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-phenyl]-propionamide;

(S)-2-Benzenesulfonylamino-N-pentyl-3-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-phenyl]-propionamide;

(S)-2-Benzenesulfonylamino-N-(4-phenyl-butyl)-3-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-phenyl]-propionamide;

(S)-2-Benzenesulfonylamino-N-(3,3-diphenyl-propyl)-3-[4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-phenyl]-propionamide;

(S)-2-Acetylamino-N-[(S)-2-[3-bromo-4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-phenyl]-1-(4-phenyl-butylcarbamoyl)-ethyl]-3-phenyl-propionamide;

(S)-2-Benzenesulfonylamino-3-[3-bromo-4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-phenyl]-N-(4-phenyl-butyl)-propionamide;

(S)-2-((S)-2-Acetylamino-3-phenyl-propionylamino)-3-[3-bromo-4-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-phenyl]-N-pentyl-propionamide; and

(S)-2-Acetylamino-N-[(S)-1-pentylcarbamoyl-2-[3-(1,1,4-trioxo-1,2,5-thiadiazolidin-2-yl)-phenyl]-ethyl]-3-phenyl-propionamide;

or a pharmaceutically acceptable salt thereof; or a prodrug derivative thereof.

17-23. (Canceled)

24. (Currently Amended) A pharmaceutical composition comprising:

a therapeutically effective amount of a compound of claim 1 in combination with one or more pharmaceutically acceptable carriers.

25-27. (Canceled)